

# Package ‘lightgbm’

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**Type** Package

**Title** Light Gradient Boosting Machine

**Version** 3.3.1

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**Description** Tree based algorithms can be improved by introducing boosting frameworks.

'LightGBM' is one such framework, based on Ke, Guolin et al. (2017) <<https://papers.nips.cc/paper/6907-lightgbm-a-highly-efficient-gradient-boosting-decision>>.

This package offers an R interface to work with it.

It is designed to be distributed and efficient with the following advantages:

1. Faster training speed and higher efficiency.
2. Lower memory usage.
3. Better accuracy.
4. Parallel learning supported.
5. Capable of handling large-scale data.

In recognition of these advantages, 'LightGBM' has been widely-used in many winning solutions of machine learning competitions.

Comparison experiments on public datasets suggest that 'LightGBM' can outperform existing boosting frameworks on both efficiency and accuracy, with significantly lower memory consumption. In addition, parallel experiments suggest that in certain circumstances, 'LightGBM' can achieve a linear speed-up in training time by using multiple machines.

**Encoding** UTF-8

**License** MIT + file LICENSE

**URL** <https://github.com/Microsoft/LightGBM>

**BugReports** <https://github.com/Microsoft/LightGBM/issues>

**NeedsCompilation** yes

**Biarch** true

**Suggests** testthat

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agaricus.test	<i>Test part from Mushroom Data Set</i>
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## Description

This data set is originally from the Mushroom data set, UCI Machine Learning Repository. This data set includes the following fields:

- label: the label for each record
- data: a sparse Matrix of dgCMatrx class, with 126 columns.

## Usage

```
data(agaricus.test)
```

## Format

A list containing a label vector, and a dgCMatrx object with 1611 rows and 126 variables

## References

<https://archive.ics.uci.edu/ml/datasets/Mushroom>

Bache, K. & Lichman, M. (2013). UCI Machine Learning Repository [<http://archive.ics.uci.edu/ml>]. Irvine, CA: University of California, School of Information and Computer Science.

---

agaricus.train      *Training part from Mushroom Data Set*

---

**Description**

This data set is originally from the Mushroom data set, UCI Machine Learning Repository. This data set includes the following fields:

- label: the label for each record
- data: a sparse Matrix of dgCMatix class, with 126 columns.

**Usage**

```
data(agaricus.train)
```

**Format**

A list containing a label vector, and a dgCMatix object with 6513 rows and 127 variables

**References**

<https://archive.ics.uci.edu/ml/datasets/Mushroom>

Bache, K. & Lichman, M. (2013). UCI Machine Learning Repository [<http://archive.ics.uci.edu/ml>]. Irvine, CA: University of California, School of Information and Computer Science.

---

bank      *Bank Marketing Data Set*

---

**Description**

This data set is originally from the Bank Marketing data set, UCI Machine Learning Repository.

It contains only the following: bank.csv with 10 randomly selected from 3 (older version of this dataset with less inputs).

**Usage**

```
data(bank)
```

**Format**

A data.table with 4521 rows and 17 variables

**References**

<http://archive.ics.uci.edu/ml/datasets/Bank+Marketing>

S. Moro, P. Cortez and P. Rita. (2014) A Data-Driven Approach to Predict the Success of Bank Telemarketing. Decision Support Systems

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dim.lgb.Dataset	<i>Dimensions of an lgb.Dataset</i>
-----------------	-------------------------------------

---

## Description

Returns a vector of numbers of rows and of columns in an lgb.Dataset.

## Usage

```
## S3 method for class 'lgb.Dataset'  
dim(x, ...)
```

## Arguments

x	Object of class lgb.Dataset
...	other parameters (ignored)

## Details

Note: since nrow and ncol internally use dim, they can also be directly used with an lgb.Dataset object.

## Value

a vector of numbers of rows and of columns

## Examples

```
data(agaricus.train, package = "lightgbm")  
train <- agaricus.train  
dtrain <- lgb.Dataset(train$data, label = train$label)
```

```
stopifnot(nrow(dtrain) == nrow(train$data))  
stopifnot(ncol(dtrain) == ncol(train$data))  
stopifnot(all(dim(dtrain) == dim(train$data)))
```

dimnames.lgb.Dataset *Handling of column names of lgb.Dataset*

---

### Description

Only column names are supported for `lgb.Dataset`, thus setting of row names would have no effect and returned row names would be `NULL`.

### Usage

```
## S3 method for class 'lgb.Dataset'
dimnames(x)

## S3 replacement method for class 'lgb.Dataset'
dimnames(x) <- value
```

### Arguments

<code>x</code>	object of class <code>lgb.Dataset</code>
<code>value</code>	a list of two elements: the first one is ignored and the second one is column names

### Details

Generic `dimnames` methods are used by `colnames`. Since row names are irrelevant, it is recommended to use `colnames` directly.

### Value

A list with the dimension names of the dataset

### Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
lgb.Dataset.construct(dtrain)
dimnames(dtrain)
colnames(dtrain)
colnames(dtrain) <- make.names(seq_len(ncol(train$data)))
print(dtrain, verbose = TRUE)
```

---

getinfo	<i>Get information of an lgb.Dataset object</i>
---------	---

---

## Description

Get one attribute of a lgb.Dataset

## Usage

```
getinfo(dataset, ...)  
  
## S3 method for class 'lgb.Dataset'  
getinfo(dataset, name, ...)
```

## Arguments

dataset	Object of class lgb.Dataset
...	other parameters (ignored)
name	the name of the information field to get (see details)

## Details

The name field can be one of the following:

- label: label lightgbm learn from ;
- weight: to do a weight rescale ;
- group: used for learning-to-rank tasks. An integer vector describing how to group rows together as ordered results from the same set of candidate results to be ranked. For example, if you have a 100-document dataset with `group = c(10, 20, 40, 10, 10, 10)`, that means that you have 6 groups, where the first 10 records are in the first group, records 11-30 are in the second group, etc.
- init\_score: initial score is the base prediction lightgbm will boost from.

## Value

info data

## Examples

```
data(agaricus.train, package = "lightgbm")  
train <- agaricus.train  
dtrain <- lgb.Dataset(train$data, label = train$label)  
lgb.Dataset.construct(dtrain)  
  
labels <- lightgbm::getinfo(dtrain, "label")  
lightgbm::setinfo(dtrain, "label", 1 - labels)
```

```
labels2 <- lightgbm::getinfo(dtrain, "label")
stopifnot(all(labels2 == 1 - labels))
```

---

get\_field

*Get one attribute of a lgb.Dataset*


---

## Description

Get one attribute of a lgb.Dataset

## Usage

```
get_field(dataset, field_name)

## S3 method for class 'lgb.Dataset'
get_field(dataset, field_name)
```

## Arguments

dataset	Object of class lgb.Dataset
field_name	String with the name of the attribute to get. One of the following. <ul style="list-style-type: none"> <li>• label: label lightgbm learns from ;</li> <li>• weight: to do a weight rescale ;</li> <li>• group: used for learning-to-rank tasks. An integer vector describing how to group rows together as ordered results from the same set of candidate results to be ranked. For example, if you have a 100-document dataset with <code>group = c(10, 20, 40, 10, 10, 10)</code>, that means that you have 6 groups, where the first 10 records are in the first group, records 11-30 are in the second group, etc.</li> <li>• <code>init_score</code>: initial score is the base prediction lightgbm will boost from.</li> </ul>

## Value

requested attribute

## Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
lgb.Dataset.construct(dtrain)

labels <- lightgbm::get_field(dtrain, "label")
lightgbm::set_field(dtrain, "label", 1 - labels)
```



```
labels2 <- lightgbm::get_field(dtrain, "label")
stopifnot(all(labels2 == 1 - labels))
```

---

`lgb.convert_with_rules`

*Data preparator for LightGBM datasets with rules (integer)*

---

## Description

Attempts to prepare a clean dataset to prepare to put in a `lgb.Dataset`. Factor, character, and logical columns are converted to integer. Missing values in factors and characters will be filled with 0L. Missing values in logicals will be filled with -1L.

This function returns and optionally takes in "rules" the describe exactly how to convert values in columns.

Columns that contain only NA values will be converted by this function but will not show up in the returned rules.

NOTE: In previous releases of LightGBM, this function was called `lgb.prepare_rules2`.

## Usage

```
lgb.convert_with_rules(data, rules = NULL)
```

## Arguments

<code>data</code>	A <code>data.frame</code> or <code>data.table</code> to prepare.
<code>rules</code>	A set of rules from the data preparator, if already used. This should be an R list, where names are column names in <code>data</code> and values are named character vectors whose names are column values and whose values are new values to replace them with.

## Value

A list with the cleaned dataset (`data`) and the rules (`rules`). Note that the data must be converted to a matrix format (`as.matrix`) for input in `lgb.Dataset`.

## Examples

```
data(iris)

str(iris)

new_iris <- lgb.convert_with_rules(data = iris)
str(new_iris$data)
```

```
data(iris) # Erase iris dataset
iris$Species[1L] <- "NEW FACTOR" # Introduce junk factor (NA)

# Use conversion using known rules
# Unknown factors become 0, excellent for sparse datasets
newer_iris <- lgb.convert_with_rules(data = iris, rules = newer_iris$rules)

# Unknown factor is now zero, perfect for sparse datasets
newer_iris$data[1L, ] # Species became 0 as it is an unknown factor

newer_iris$data[1L, 5L] <- 1.0 # Put back real initial value

# Is the newly created dataset equal? YES!
all.equal(newer_iris$data, newer_iris$data)

# Can we test our own rules?
data(iris) # Erase iris dataset

# We remapped values differently
personal_rules <- list(
  Species = c(
    "setosa" = 3L
    , "versicolor" = 2L
    , "virginica" = 1L
  )
)
newest_iris <- lgb.convert_with_rules(data = iris, rules = personal_rules)
str(newest_iris$data) # SUCCESS!
```

---

lgb.cv

*Main CV logic for LightGBM*

---

## Description

Cross validation logic used by LightGBM

## Usage

```
lgb.cv(
  params = list(),
  data,
  nrounds = 100L,
  nfold = 3L,
  label = NULL,
  weight = NULL,
  obj = NULL,
  eval = NULL,
  verbose = 1L,
```

```

record = TRUE,
eval_freq = 1L,
showsd = TRUE,
stratified = TRUE,
folds = NULL,
init_model = NULL,
colnames = NULL,
categorical_feature = NULL,
early_stopping_rounds = NULL,
callbacks = list(),
reset_data = FALSE,
...
)

```

## Arguments

params	a list of parameters. See the <a href="#">"Parameters" section of the documentation</a> for a list of parameters and valid values.
data	a <code>lgb.Dataset</code> object, used for training. Some functions, such as <code>lgb.cv</code> , may allow you to pass other types of data like <code>matrix</code> and then separately supply <code>label</code> as a keyword argument.
nrounds	number of training rounds
nfold	the original dataset is randomly partitioned into <code>nfold</code> equal size subsamples.
label	Vector of labels, used if data is not an <code>lgb.Dataset</code>
weight	vector of response values. If not <code>NULL</code> , will set to dataset
obj	objective function, can be character or custom objective function. Examples include <code>regression</code> , <code>regression_l1</code> , <code>huber</code> , <code>binary</code> , <code>lambda_rank</code> , <code>multiclass</code> , <code>multiclass</code>
eval	evaluation function(s). This can be a character vector, function, or list with a mixture of strings and functions. <ul style="list-style-type: none"> <li>• <b>a. character vector:</b> If you provide a character vector to this argument, it should contain strings with valid evaluation metrics. See <a href="#">The "metric" section of the documentation</a> for a list of valid metrics.</li> <li>• <b>b. function:</b> You can provide a custom evaluation function. This should accept the keyword arguments <code>preds</code> and <code>dtrain</code> and should return a named list with three elements: <ul style="list-style-type: none"> <li>– <code>name</code>: A string with the name of the metric, used for printing and storing results.</li> <li>– <code>value</code>: A single number indicating the value of the metric for the given predictions and true values</li> <li>– <code>higher_better</code>: A boolean indicating whether higher values indicate a better fit. For example, this would be <code>FALSE</code> for metrics like MAE or RMSE.</li> </ul> </li> <li>• <b>c. list:</b> If a list is given, it should only contain character vectors and functions. These should follow the requirements from the descriptions above.</li> </ul>

verbose	verbosity for output, if $\leq 0$ , also will disable the print of evaluation during training
record	Boolean, TRUE will record iteration message to <code>booster\$record_evals</code>
eval_freq	evaluation output frequency, only effect when <code>verbose &gt; 0</code>
showsd	boolean, whether to show standard deviation of cross validation. This parameter defaults to TRUE. Setting it to FALSE can lead to a slight speedup by avoiding unnecessary computation.
stratified	a boolean indicating whether sampling of folds should be stratified by the values of outcome labels.
folds	list provides a possibility to use a list of pre-defined CV folds (each element must be a vector of test fold's indices). When folds are supplied, the <code>ifold</code> and <code>stratified</code> parameters are ignored.
init_model	path of model file of <code>lgb.Booster</code> object, will continue training from this model
colnames	feature names, if not null, will use this to overwrite the names in dataset
categorical_feature	categorical features. This can either be a character vector of feature names or an integer vector with the indices of the features (e.g. <code>c(1L, 10L)</code> to say "the first and tenth columns").
early_stopping_rounds	int. Activates early stopping. When this parameter is non-null, training will stop if the evaluation of any metric on any validation set fails to improve for <code>early_stopping_rounds</code> consecutive boosting rounds. If training stops early, the returned model will have attribute <code>best_iter</code> set to the iteration number of the best iteration.
callbacks	List of callback functions that are applied at each iteration.
reset_data	Boolean, setting it to TRUE (not the default value) will transform the booster model into a predictor model which frees up memory and the original datasets
...	other parameters, see <code>Parameters.rst</code> for more information. A few key parameters: <ul style="list-style-type: none"> <li>• <code>boosting</code>: Boosting type. "gbdt", "rf", "dart" or "goss".</li> <li>• <code>num_leaves</code>: Maximum number of leaves in one tree.</li> <li>• <code>max_depth</code>: Limit the max depth for tree model. This is used to deal with overfit when #data is small. Tree still grow by leaf-wise.</li> <li>• <code>num_threads</code>: Number of threads for LightGBM. For the best speed, set this to the number of real CPU cores(<code>parallel::detectCores(logical = FALSE)</code>), not the number of threads (most CPU using hyper-threading to generate 2 threads per CPU core).</li> </ul> <p>NOTE: As of v3.3.0, use of ... is deprecated. Add parameters to <code>params</code> directly.</p>

### Value

a trained model `lgb.CVBooster`.

## Early Stopping

"early stopping" refers to stopping the training process if the model's performance on a given validation set does not improve for several consecutive iterations.

If multiple arguments are given to `eval`, their order will be preserved. If you enable early stopping by setting `early_stopping_rounds` in `params`, by default all metrics will be considered for early stopping.

If you want to only consider the first metric for early stopping, pass `first_metric_only = TRUE` in `params`. Note that if you also specify `metric` in `params`, that metric will be considered the "first" one. If you omit `metric`, a default metric will be used based on your choice for the parameter `obj` (keyword argument) or `objective` (passed into `params`).

## Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
model <- lgb.cv(
  params = params
  , data = dtrain
  , nrounds = 5L
  , nfold = 3L
)
```

---

lgb.Dataset

*Construct lgb.Dataset object*


---

## Description

Construct `lgb.Dataset` object from dense matrix, sparse matrix or local file (that was created previously by saving an `lgb.Dataset`).

## Usage

```
lgb.Dataset(
  data,
  params = list(),
  reference = NULL,
  colnames = NULL,
  categorical_feature = NULL,
```

```

    free_raw_data = TRUE,
    info = list(),
    label = NULL,
    weight = NULL,
    group = NULL,
    init_score = NULL,
    ...
)

```

### Arguments

data	a matrix object, a dgCMatrix object, a character representing a path to a text file (CSV, TSV, or LibSVM), or a character representing a path to a binary lgb.Dataset file
params	a list of parameters. See <a href="#">The "Dataset Parameters" section of the documentation</a> for a list of parameters and valid values.
reference	reference dataset. When LightGBM creates a Dataset, it does some preprocessing like binning continuous features into histograms. If you want to apply the same bin boundaries from an existing dataset to new data, pass that existing Dataset to this argument.
colnames	names of columns
categorical_feature	categorical features. This can either be a character vector of feature names or an integer vector with the indices of the features (e.g. <code>c(1L, 10L)</code> to say "the first and tenth columns").
free_raw_data	LightGBM constructs its data format, called a "Dataset", from tabular data. By default, that Dataset object on the R side does not keep a copy of the raw data. This reduces LightGBM's memory consumption, but it means that the Dataset object cannot be changed after it has been constructed. If you'd prefer to be able to change the Dataset object after construction, set <code>free_raw_data = FALSE</code> .
info	a list of information of the lgb.Dataset object. NOTE: use of info is deprecated as of v3.3.0. Use keyword arguments (e.g. <code>init_score = init_score</code> ) directly.
label	vector of labels to use as the target variable
weight	numeric vector of sample weights
group	used for learning-to-rank tasks. An integer vector describing how to group rows together as ordered results from the same set of candidate results to be ranked. For example, if you have a 100-document dataset with <code>group = c(10, 20, 40, 10, 10, 10)</code> , that means that you have 6 groups, where the first 10 records are in the first group, records 11-30 are in the second group, etc.
init_score	initial score is the base prediction lightgbm will boost from
...	other parameters passed to params

### Value

constructed dataset

## Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data_file <- tempfile(fileext = ".data")
lgb.Dataset.save(dtrain, data_file)
dtrain <- lgb.Dataset(data_file)
lgb.Dataset.construct(dtrain)
```

---

`lgb.Dataset.construct` *Construct Dataset explicitly*

---

## Description

Construct Dataset explicitly

## Usage

```
lgb.Dataset.construct(dataset)
```

## Arguments

dataset            Object of class `lgb.Dataset`

## Value

constructed dataset

## Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
lgb.Dataset.construct(dtrain)
```

---

```
lgb.Dataset.create.valid
```

*Construct validation data*

---

## Description

Construct validation data according to training data

## Usage

```
lgb.Dataset.create.valid(
  dataset,
  data,
  info = list(),
  label = NULL,
  weight = NULL,
  group = NULL,
  init_score = NULL,
  params = list(),
  ...
)
```

## Arguments

dataset	lgb.Dataset object, training data
data	a matrix object, a dgCMatix object, a character representing a path to a text file (CSV, TSV, or LibSVM), or a character representing a path to a binary Dataset file
info	a list of information of the lgb.Dataset object. NOTE: use of info is deprecated as of v3.3.0. Use keyword arguments (e.g. <code>init_score = init_score</code> ) directly.
label	vector of labels to use as the target variable
weight	numeric vector of sample weights
group	used for learning-to-rank tasks. An integer vector describing how to group rows together as ordered results from the same set of candidate results to be ranked. For example, if you have a 100-document dataset with <code>group = c(10, 20, 40, 10, 10, 10)</code> , that means that you have 6 groups, where the first 10 records are in the first group, records 11-30 are in the second group, etc.
init_score	initial score is the base prediction lightgbm will boost from
params	a list of parameters. See <a href="#">The "Dataset Parameters" section of the documentation</a> for a list of parameters and valid values. If this is an empty list (the default), the validation Dataset will have the same parameters as the Dataset passed to argument dataset.
...	additional lgb.Dataset parameters. NOTE: As of v3.3.0, use of ... is deprecated. Add parameters to params directly.



**Value**

constructed dataset

**Examples**

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)

# parameters can be changed between the training data and validation set,
# for example to account for training data in a text file with a header row
# and validation data in a text file without it
train_file <- tempfile(pattern = "train_", fileext = ".csv")
write.table(
  data.frame(y = rnorm(100L), x1 = rnorm(100L), x2 = rnorm(100L))
  , file = train_file
  , sep = ","
  , col.names = TRUE
  , row.names = FALSE
  , quote = FALSE
)

valid_file <- tempfile(pattern = "valid_", fileext = ".csv")
write.table(
  data.frame(y = rnorm(100L), x1 = rnorm(100L), x2 = rnorm(100L))
  , file = valid_file
  , sep = ","
  , col.names = FALSE
  , row.names = FALSE
  , quote = FALSE
)

dtrain <- lgb.Dataset(
  data = train_file
  , params = list(has_header = TRUE)
)
dtrain$construct()

dvalid <- lgb.Dataset(
  data = valid_file
  , params = list(has_header = FALSE)
)
dvalid$construct()
```

`lgb.Dataset.save`      *Save lgb.Dataset to a binary file*

---

### Description

Please note that `init_score` is not saved in binary file. If you need it, please set it again after loading Dataset.

### Usage

```
lgb.Dataset.save(dataset, fname)
```

### Arguments

<code>dataset</code>	object of class <code>lgb.Dataset</code>
<code>fname</code>	object filename of output file

### Value

the dataset you passed in

### Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
lgb.Dataset.save(dtrain, tempfile(fileext = ".bin"))
```

---

`lgb.Dataset.set.categorical`  
*Set categorical feature of lgb.Dataset*

---

### Description

Set the categorical features of an `lgb.Dataset` object. Use this function to tell LightGBM which features should be treated as categorical.

### Usage

```
lgb.Dataset.set.categorical(dataset, categorical_feature)
```

**Arguments**

`dataset` object of class `lgb.Dataset`  
`categorical_feature` categorical features. This can either be a character vector of feature names or an integer vector with the indices of the features (e.g. `c(1L, 10L)` to say "the first and tenth columns").

**Value**

the dataset you passed in

**Examples**

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data_file <- tempfile(fileext = ".data")
lgb.Dataset.save(dtrain, data_file)
dtrain <- lgb.Dataset(data_file)
lgb.Dataset.set.categorical(dtrain, 1L:2L)
```

---

`lgb.Dataset.set.reference`  
*Set reference of lgb.Dataset*

---

**Description**

If you want to use validation data, you should set reference to training data

**Usage**

```
lgb.Dataset.set.reference(dataset, reference)
```

**Arguments**

`dataset` object of class `lgb.Dataset`  
`reference` object of class `lgb.Dataset`

**Value**

the dataset you passed in

## Examples

```
# create training Dataset
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)

# create a validation Dataset, using dtrain as a reference
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset(test$data, label = test$label)
lgb.Dataset.set.reference(dtest, dtrain)
```

---

lgb.dump

*Dump LightGBM model to json*

---

## Description

Dump LightGBM model to json

## Usage

```
lgb.dump(booster, num_iteration = NULL)
```

## Arguments

**booster**            Object of class `lgb.Booster`  
**num\_iteration**    number of iteration want to predict with, `NULL` or `<= 0` means use best iteration

## Value

json format of model

## Examples

```
library(lightgbm)
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
```

```
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 10L
  , valids = valids
  , early_stopping_rounds = 5L
)
json_model <- lgb.dump(model)
```

---

`lgb.get.eval.result`    *Get record evaluation result from booster*

---

## Description

Given a `lgb.Booster`, return evaluation results for a particular metric on a particular dataset.

## Usage

```
lgb.get.eval.result(
  booster,
  data_name,
  eval_name,
  iters = NULL,
  is_err = FALSE
)
```

## Arguments

<code>booster</code>	Object of class <code>lgb.Booster</code>
<code>data_name</code>	Name of the dataset to return evaluation results for.
<code>eval_name</code>	Name of the evaluation metric to return results for.
<code>iters</code>	An integer vector of iterations you want to get evaluation results for. If <code>NULL</code> (the default), evaluation results for all iterations will be returned.
<code>is_err</code>	<code>TRUE</code> will return evaluation error instead

## Value

numeric vector of evaluation result

**Examples**

```

# train a regression model
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 5L
  , valids = valids
)

# Examine valid data_name values
print(setdiff(names(model$record_evals), "start_iter"))

# Examine valid eval_name values for dataset "test"
print(names(model$record_evals[["test"]]))

# Get L2 values for "test" dataset
lgb.get.eval.result(model, "test", "l2")

```

---

lgb.importance

*Compute feature importance in a model*


---

**Description**

Creates a data.table of feature importances in a model.

**Usage**

```
lgb.importance(model, percentage = TRUE)
```

**Arguments**

model	object of class lgb.Booster.
percentage	whether to show importance in relative percentage.

**Value**

For a tree model, a data.table with the following columns:

- Feature: Feature names in the model.
- Gain: The total gain of this feature's splits.
- Cover: The number of observation related to this feature.
- Frequency: The number of times a feature splitted in trees.

**Examples**

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)

params <- list(
  objective = "binary"
  , learning_rate = 0.1
  , max_depth = -1L
  , min_data_in_leaf = 1L
  , min_sum_hessian_in_leaf = 1.0
)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 5L
)

tree_imp1 <- lgb.importance(model, percentage = TRUE)
tree_imp2 <- lgb.importance(model, percentage = FALSE)
```

---

lgb.interprete

*Compute feature contribution of prediction*


---

**Description**

Computes feature contribution components of rawscore prediction.

**Usage**

```
lgb.interprete(model, data, idxset, num_iteration = NULL)
```

**Arguments**

model	object of class lgb.Booster.
data	a matrix object or a dgCMatix object.
idxset	an integer vector of indices of rows needed.
num_iteration	number of iteration want to predict with, NULL or <= 0 means use best iteration.

**Value**

For regression, binary classification and lambdarank model, a list of `data.table` with the following columns:

- **Feature:** Feature names in the model.
- **Contribution:** The total contribution of this feature's splits.

For multiclass classification, a list of `data.table` with the `Feature` column and `Contribution` columns to each class.

**Examples**

```
Logit <- function(x) log(x / (1.0 - x))
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
set_field(
  dataset = dtrain
  , field_name = "init_score"
  , data = rep(Logit(mean(train$label)), length(train$label))
)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test

params <- list(
  objective = "binary"
  , learning_rate = 0.1
  , max_depth = -1L
  , min_data_in_leaf = 1L
  , min_sum_hessian_in_leaf = 1.0
)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 3L
)

tree_interpretation <- lgb.interprete(model, test$data, 1L:5L)
```

---

lgb.load

*Load LightGBM model*


---

**Description**

Load LightGBM takes in either a file path or model string. If both are provided, Load will default to loading from file



**Usage**

```
lgb.load(filename = NULL, model_str = NULL)
```

**Arguments**

filename	path of model file
model_str	a str containing the model

**Value**

lgb.Booster

**Examples**

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 5L
  , valids = valids
  , early_stopping_rounds = 3L
)
model_file <- tempfile(fileext = ".txt")
lgb.save(model, model_file)
load_booster <- lgb.load(filename = model_file)
model_string <- model$save_model_to_string(NULL) # saves best iteration
load_booster_from_str <- lgb.load(model_str = model_string)
```

---

lgb.model.dt.tree      *Parse a LightGBM model json dump*

---

**Description**

Parse a LightGBM model json dump into a data.table structure.

**Usage**

```
lgb.model.dt.tree(model, num_iteration = NULL)
```

**Arguments**

```
model          object of class lgb.Booster
num_iteration  number of iterations you want to predict with. NULL or <= 0 means use best
               iteration
```

**Value**

A data.table with detailed information about model trees' nodes and leaves.

The columns of the data.table are:

- tree\_index: ID of a tree in a model (integer)
- split\_index: ID of a node in a tree (integer)
- split\_feature: for a node, it's a feature name (character); for a leaf, it simply labels it as "NA"
- node\_parent: ID of the parent node for current node (integer)
- leaf\_index: ID of a leaf in a tree (integer)
- leaf\_parent: ID of the parent node for current leaf (integer)
- split\_gain: Split gain of a node
- threshold: Splitting threshold value of a node
- decision\_type: Decision type of a node
- default\_left: Determine how to handle NA value, TRUE -> Left, FALSE -> Right
- internal\_value: Node value
- internal\_count: The number of observation collected by a node
- leaf\_value: Leaf value
- leaf\_count: The number of observation collected by a leaf

**Examples**

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)

params <- list(
  objective = "binary"
  , learning_rate = 0.01
  , num_leaves = 63L
  , max_depth = -1L
  , min_data_in_leaf = 1L
  , min_sum_hessian_in_leaf = 1.0
)
```

```
model <- lgb.train(params, dtrain, 10L)
tree_dt <- lgb.model.dt.tree(model)
```

---

`lgb.plot.importance` *Plot feature importance as a bar graph*

---

### Description

Plot previously calculated feature importance: Gain, Cover and Frequency, as a bar graph.

### Usage

```
lgb.plot.importance(
  tree_imp,
  top_n = 10L,
  measure = "Gain",
  left_margin = 10L,
  cex = NULL
)
```

### Arguments

<code>tree_imp</code>	a data.table returned by <a href="#">lgb.importance</a> .
<code>top_n</code>	maximal number of top features to include into the plot.
<code>measure</code>	the name of importance measure to plot, can be "Gain", "Cover" or "Frequency".
<code>left_margin</code>	(base R barplot) allows to adjust the left margin size to fit feature names.
<code>cex</code>	(base R barplot) passed as <code>cex.names</code> parameter to <a href="#">barplot</a> . Set a number smaller than 1.0 to make the bar labels smaller than R's default and values greater than 1.0 to make them larger.

### Details

The graph represents each feature as a horizontal bar of length proportional to the defined importance of a feature. Features are shown ranked in a decreasing importance order.

### Value

The `lgb.plot.importance` function creates a barplot and silently returns a processed data.table with `top_n` features sorted by defined importance.

**Examples**

```

data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)

params <- list(
  objective = "binary"
  , learning_rate = 0.1
  , min_data_in_leaf = 1L
  , min_sum_hessian_in_leaf = 1.0
)

model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 5L
)

tree_imp <- lgb.importance(model, percentage = TRUE)
lgb.plot.importance(tree_imp, top_n = 5L, measure = "Gain")

```

---

lgb.plot.interpretation

*Plot feature contribution as a bar graph*


---

**Description**

Plot previously calculated feature contribution as a bar graph.

**Usage**

```

lgb.plot.interpretation(
  tree_interpretation_dt,
  top_n = 10L,
  cols = 1L,
  left_margin = 10L,
  cex = NULL
)

```

**Arguments**

tree_interpretation_dt	a data.table returned by <a href="#">lgb.interprete</a> .
top_n	maximal number of top features to include into the plot.
cols	the column numbers of layout, will be used only for multiclass classification feature contribution.

`left_margin` (base R barplot) allows to adjust the left margin size to fit feature names.  
`cex` (base R barplot) passed as `cex.names` parameter to `barplot`.

### Details

The graph represents each feature as a horizontal bar of length proportional to the defined contribution of a feature. Features are shown ranked in a decreasing contribution order.

### Value

The `lgb.plot.interpretation` function creates a barplot.

### Examples

```
Logit <- function(x) {
  log(x / (1.0 - x))
}
data(agaricus.train, package = "lightgbm")
labels <- agaricus.train$label
dtrain <- lgb.Dataset(
  agaricus.train$data
  , label = labels
)
set_field(
  dataset = dtrain
  , field_name = "init_score"
  , data = rep(Logit(mean(labels)), length(labels))
)

data(agaricus.test, package = "lightgbm")

params <- list(
  objective = "binary"
  , learning_rate = 0.1
  , max_depth = -1L
  , min_data_in_leaf = 1L
  , min_sum_hessian_in_leaf = 1.0
)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 5L
)

tree_interpretation <- lgb.interprete(
  model = model
  , data = agaricus.test$data
  , idxset = 1L:5L
)
lgb.plot.interpretation(
  tree_interpretation_dt = tree_interpretation[[1L]]
```

```

    , top_n = 3L
  )

```

---

lgb.save

*Save LightGBM model*


---

## Description

Save LightGBM model

## Usage

```
lgb.save(booster, filename, num_iteration = NULL)
```

## Arguments

booster	Object of class lgb.Booster
filename	saved filename
num_iteration	number of iteration want to predict with, NULL or <= 0 means use best iteration

## Value

lgb.Booster

## Examples

```

library(lightgbm)
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 10L
  , valids = valids
  , early_stopping_rounds = 5L
)

```

```
lgb.save(model, tempfile(fileext = ".txt"))
```

---

lgb.train

*Main training logic for LightGBM*


---

## Description

Logic to train with LightGBM

## Usage

```
lgb.train(
  params = list(),
  data,
  nrounds = 100L,
  valids = list(),
  obj = NULL,
  eval = NULL,
  verbose = 1L,
  record = TRUE,
  eval_freq = 1L,
  init_model = NULL,
  colnames = NULL,
  categorical_feature = NULL,
  early_stopping_rounds = NULL,
  callbacks = list(),
  reset_data = FALSE,
  ...
)
```

## Arguments

params	a list of parameters. See <a href="#">the "Parameters" section of the documentation</a> for a list of parameters and valid values.
data	a <code>lgb.Dataset</code> object, used for training. Some functions, such as <code>lgb.cv</code> , may allow you to pass other types of data like matrix and then separately supply label as a keyword argument.
nrounds	number of training rounds
valids	a list of <code>lgb.Dataset</code> objects, used for validation
obj	objective function, can be character or custom objective function. Examples include regression, regression_l1, huber, binary, lambdarank, multiclass, multiclass
eval	evaluation function(s). This can be a character vector, function, or list with a mixture of strings and functions.

- **a. character vector:** If you provide a character vector to this argument, it should contain strings with valid evaluation metrics. See [The "metric" section of the documentation](#) for a list of valid metrics.
- **b. function:** You can provide a custom evaluation function. This should accept the keyword arguments `preds` and `dtrain` and should return a named list with three elements:
  - `name`: A string with the name of the metric, used for printing and storing results.
  - `value`: A single number indicating the value of the metric for the given predictions and true values
  - `higher_better`: A boolean indicating whether higher values indicate a better fit. For example, this would be `FALSE` for metrics like MAE or RMSE.
- **c. list:** If a list is given, it should only contain character vectors and functions. These should follow the requirements from the descriptions above.

<code>verbose</code>	verbosity for output, if $\leq 0$ , also will disable the print of evaluation during training
<code>record</code>	Boolean, <code>TRUE</code> will record iteration message to <code>booster\$record_evals</code>
<code>eval_freq</code>	evaluation output frequency, only effect when <code>verbose &gt; 0</code>
<code>init_model</code>	path of model file of <code>lgb</code> . Booster object, will continue training from this model
<code>colnames</code>	feature names, if not null, will use this to overwrite the names in dataset
<code>categorical_feature</code>	categorical features. This can either be a character vector of feature names or an integer vector with the indices of the features (e.g. <code>c(1L, 10L)</code> to say "the first and tenth columns").
<code>early_stopping_rounds</code>	int. Activates early stopping. When this parameter is non-null, training will stop if the evaluation of any metric on any validation set fails to improve for <code>early_stopping_rounds</code> consecutive boosting rounds. If training stops early, the returned model will have attribute <code>best_iter</code> set to the iteration number of the best iteration.
<code>callbacks</code>	List of callback functions that are applied at each iteration.
<code>reset_data</code>	Boolean, setting it to <code>TRUE</code> (not the default value) will transform the booster model into a predictor model which frees up memory and the original datasets
<code>...</code>	other parameters, see <a href="#">the "Parameters" section of the documentation</a> for more information. A few key parameters: <ul style="list-style-type: none"> <li>• <code>boosting</code>: Boosting type. "gbdt", "rf", "dart" or "goss".</li> <li>• <code>num_leaves</code>: Maximum number of leaves in one tree.</li> <li>• <code>max_depth</code>: Limit the max depth for tree model. This is used to deal with overfitting. Tree still grow by leaf-wise.</li> <li>• <code>num_threads</code>: Number of threads for LightGBM. For the best speed, set this to the number of real CPU cores(<code>parallel::detectCores(logical = FALSE)</code>), not the number of threads (most CPU using hyper-threading to generate 2 threads per CPU core).</li> </ul>

NOTE: As of v3.3.0, use of `...` is deprecated. Add parameters to `params` directly.



**Value**

a trained booster model `lgb.Booster`.

**Early Stopping**

"early stopping" refers to stopping the training process if the model's performance on a given validation set does not improve for several consecutive iterations.

If multiple arguments are given to `eval`, their order will be preserved. If you enable early stopping by setting `early_stopping_rounds` in `params`, by default all metrics will be considered for early stopping.

If you want to only consider the first metric for early stopping, pass `first_metric_only = TRUE` in `params`. Note that if you also specify `metric` in `params`, that metric will be considered the "first" one. If you omit `metric`, a default metric will be used based on your choice for the parameter `obj` (keyword argument) or `objective` (passed into `params`).

**Examples**

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 5L
  , valids = valids
  , early_stopping_rounds = 3L
)
```

---

lgb.unloader

*Remove lightgbm and its objects from an environment*


---

**Description**

Attempts to unload LightGBM packages so you can remove objects cleanly without having to restart R. This is useful for instance if an object becomes stuck for no apparent reason and you do not want to restart R to fix the lost object.

**Usage**

```
lgb.unloader(restore = TRUE, wipe = FALSE, envir = .GlobalEnv)
```

**Arguments**

restore	Whether to reload LightGBM immediately after detaching from R. Defaults to TRUE which means automatically reload LightGBM once unloading is performed.
wipe	Whether to wipe all lgb.Dataset and lgb.Booster from the global environment. Defaults to FALSE which means to not remove them.
envir	The environment to perform wiping on if wipe == TRUE. Defaults to .GlobalEnv which is the global environment.

**Value**

NULL invisibly.

**Examples**

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 5L
  , valids = valids
)

lgb.unloader(restore = FALSE, wipe = FALSE, envir = .GlobalEnv)
rm(model, dtrain, dtest) # Not needed if wipe = TRUE
gc() # Not needed if wipe = TRUE

library(lightgbm)
# Do whatever you want again with LightGBM without object clashing
```

lightgbm

*Train a LightGBM model***Description**

Simple interface for training a LightGBM model.

**Usage**

```
lightgbm(
  data,
  label = NULL,
  weight = NULL,
  params = list(),
  nrounds = 100L,
  verbose = 1L,
  eval_freq = 1L,
  early_stopping_rounds = NULL,
  save_name = "lightgbm.model",
  init_model = NULL,
  callbacks = list(),
  ...
)
```

**Arguments**

data	a <code>lgb.Dataset</code> object, used for training. Some functions, such as <code>lgb.cv</code> , may allow you to pass other types of data like matrix and then separately supply label as a keyword argument.
label	Vector of labels, used if data is not an <code>lgb.Dataset</code>
weight	vector of response values. If not <code>NULL</code> , will set to dataset
params	a list of parameters. See the " <b>Parameters</b> " section of the documentation for a list of parameters and valid values.
nrounds	number of training rounds
verbose	verbosity for output, if $\leq 0$ , also will disable the print of evaluation during training
eval_freq	evaluation output frequency, only effect when <code>verbose &gt; 0</code>
early_stopping_rounds	int. Activates early stopping. When this parameter is non-null, training will stop if the evaluation of any metric on any validation set fails to improve for <code>early_stopping_rounds</code> consecutive boosting rounds. If training stops early, the returned model will have attribute <code>best_iter</code> set to the iteration number of the best iteration.
save_name	File name to use when writing the trained model to disk. Should end in ".model".

<code>init_model</code>	path of model file of <code>lgb.Booster</code> object, will continue training from this model
<code>callbacks</code>	List of callback functions that are applied at each iteration.
<code>...</code>	Additional arguments passed to <code>lgb.train</code> . For example <ul style="list-style-type: none"> <li>• <code>valids</code>: a list of <code>lgb.Dataset</code> objects, used for validation</li> <li>• <code>obj</code>: objective function, can be character or custom objective function. Examples include <code>regression</code>, <code>regression_l1</code>, <code>huber</code>, <code>binary</code>, <code>lambdarank</code>, <code>multiclass</code>, <code>multiclass</code></li> <li>• <code>eval</code>: evaluation function, can be (a list of) character or custom eval function</li> <li>• <code>record</code>: Boolean, TRUE will record iteration message to <code>booster\$record_evals</code></li> <li>• <code>colnames</code>: feature names, if not null, will use this to overwrite the names in dataset</li> <li>• <code>categorical_feature</code>: categorical features. This can either be a character vector of feature names or an integer vector with the indices of the features (e.g. <code>c(1L, 10L)</code> to say "the first and tenth columns").</li> <li>• <code>reset_data</code>: Boolean, setting it to TRUE (not the default value) will transform the booster model into a predictor model which frees up memory and the original datasets</li> <li>• <code>boosting</code>: Boosting type. "gbdt", "rf", "dart" or "goss".</li> <li>• <code>num_leaves</code>: Maximum number of leaves in one tree.</li> <li>• <code>max_depth</code>: Limit the max depth for tree model. This is used to deal with overfit when #data is small. Tree still grow by leaf-wise.</li> <li>• <code>num_threads</code>: Number of threads for LightGBM. For the best speed, set this to the number of real CPU cores(<code>parallel::detectCores(logical = FALSE)</code>), not the number of threads (most CPU using hyper-threading to generate 2 threads per CPU core).</li> </ul>

## Value

a trained `lgb.Booster`

## Early Stopping

"early stopping" refers to stopping the training process if the model's performance on a given validation set does not improve for several consecutive iterations.

If multiple arguments are given to `eval`, their order will be preserved. If you enable early stopping by setting `early_stopping_rounds` in `params`, by default all metrics will be considered for early stopping.

If you want to only consider the first metric for early stopping, pass `first_metric_only = TRUE` in `params`. Note that if you also specify `metric` in `params`, that metric will be considered the "first" one. If you omit `metric`, a default metric will be used based on your choice for the parameter `obj` (keyword argument) or `objective` (passed into `params`).

---

predict.lgb.Booster     *Predict method for LightGBM model*

---

### Description

Predicted values based on class lgb.Booster

### Usage

```
## S3 method for class 'lgb.Booster'
predict(
  object,
  data,
  start_iteration = NULL,
  num_iteration = NULL,
  rawscore = FALSE,
  predleaf = FALSE,
  predcontrib = FALSE,
  header = FALSE,
  reshape = FALSE,
  params = list(),
  ...
)
```

### Arguments

object	Object of class lgb.Booster
data	a matrix object, a dgCMatix object or a character representing a path to a text file (CSV, TSV, or LibSVM)
start_iteration	int or None, optional (default=None) Start index of the iteration to predict. If None or <= 0, starts from the first iteration.
num_iteration	int or None, optional (default=None) Limit number of iterations in the prediction. If None, if the best iteration exists and start_iteration is None or <= 0, the best iteration is used; otherwise, all iterations from start_iteration are used. If <= 0, all iterations from start_iteration are used (no limits).
rawscore	whether the prediction should be returned in the form of original untransformed sum of predictions from boosting iterations' results. E.g., setting rawscore=TRUE for logistic regression would result in predictions for log-odds instead of probabilities.
predleaf	whether predict leaf index instead.
predcontrib	return per-feature contributions for each record.
header	only used for prediction for text file. True if text file has header
reshape	whether to reshape the vector of predictions to a matrix form when there are several prediction outputs per case.

`params` a list of additional named parameters. See the ["Predict Parameters" section of the documentation](#) for a list of parameters and valid values.

`...` Additional prediction parameters. NOTE: deprecated as of v3.3.0. Use `params` instead.

## Value

For regression or binary classification, it returns a vector of length `nrows(data)`. For multiclass classification, either a `num_class * nrows(data)` vector or a `(nrows(data), num_class)` dimension matrix is returned, depending on the `reshape` value.

When `predleaf = TRUE`, the output is a matrix object with the number of columns corresponding to the number of trees.

## Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 5L
  , valids = valids
)
preds <- predict(model, test$data)

# pass other prediction parameters
preds <- predict(
  model,
  test$data,
  params = list(
    predict_disable_shape_check = TRUE
  )
)
```

---

readRDS.lgb.Booster    *readRDS for lgb.Booster models*

---

## Description

Attempts to load a model stored in a .rds file, using [readRDS](#)

## Usage

```
readRDS.lgb.Booster(file, refhook = NULL)
```

## Arguments

**file**                    a connection or the name of the file where the R object is saved to or read from.  
**refhook**                a hook function for handling reference objects.

## Value

lgb.Booster

## Examples

```
library(lightgbm)
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 10L
  , valids = valids
  , early_stopping_rounds = 5L
)
model_file <- tempfile(fileext = ".rds")
saveRDS.lgb.Booster(model, model_file)
new_model <- readRDS.lgb.Booster(model_file)
```

---

saveRDS.lgb.Booster     *saveRDS for lgb.Booster models*

---

### Description

Attempts to save a model using RDS. Has an additional parameter (`raw`) which decides whether to save the raw model or not.

### Usage

```
saveRDS.lgb.Booster(  
  object,  
  file,  
  ascii = FALSE,  
  version = NULL,  
  compress = TRUE,  
  refhook = NULL,  
  raw = TRUE  
)
```

### Arguments

<code>object</code>	<code>lgb.Booster</code> object to serialize.
<code>file</code>	a connection or the name of the file where the R object is saved to or read from.
<code>ascii</code>	a logical. If <code>TRUE</code> or <code>NA</code> , an ASCII representation is written; otherwise (default), a binary one is used. See the comments in the help for <code>save</code> .
<code>version</code>	the workspace format version to use. <code>NULL</code> specifies the current default version (2). Versions prior to 2 are not supported, so this will only be relevant when there are later versions.
<code>compress</code>	a logical specifying whether saving to a named file is to use "gzip" compression, or one of "gzip", "bzip2" or "xz" to indicate the type of compression to be used. Ignored if file is a connection.
<code>refhook</code>	a hook function for handling reference objects.
<code>raw</code>	whether to save the model in a raw variable or not, recommended to leave it to <code>TRUE</code> .

### Value

`NULL` invisibly.

### Examples

```
library(lightgbm)  
data(agaricus.train, package = "lightgbm")  
train <- agaricus.train
```



```

dtrain <- lgb.Dataset(train$data, label = train$label)
data(agaricus.test, package = "lightgbm")
test <- agaricus.test
dtest <- lgb.Dataset.create.valid(dtrain, test$data, label = test$label)
params <- list(
  objective = "regression"
  , metric = "l2"
  , min_data = 1L
  , learning_rate = 1.0
)
valids <- list(test = dtest)
model <- lgb.train(
  params = params
  , data = dtrain
  , nrounds = 10L
  , valids = valids
  , early_stopping_rounds = 5L
)
model_file <- tempfile(fileext = ".rds")
saveRDS.lgb.Booster(model, model_file)

```

---

setinfo

*Set information of an lgb.Dataset object*


---

## Description

Set one attribute of a lgb.Dataset

## Usage

```
setinfo(dataset, ...)
```

```
## S3 method for class 'lgb.Dataset'
setinfo(dataset, name, info, ...)
```

## Arguments

dataset	Object of class lgb.Dataset
...	other parameters (ignored)
name	the name of the field to get
info	the specific field of information to set

## Details

The name field can be one of the following:

- label: vector of labels to use as the target variable

- `weight`: to do a weight rescale
- `init_score`: initial score is the base prediction lightgbm will boost from
- `group`: used for learning-to-rank tasks. An integer vector describing how to group rows together as ordered results from the same set of candidate results to be ranked. For example, if you have a 100-document dataset with `group = c(10, 20, 40, 10, 10, 10)`, that means that you have 6 groups, where the first 10 records are in the first group, records 11-30 are in the second group, etc.

### Value

the dataset you passed in

### Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
lgb.Dataset.construct(dtrain)

labels <- lightgbm::getinfo(dtrain, "label")
lightgbm::setinfo(dtrain, "label", 1 - labels)

labels2 <- lightgbm::getinfo(dtrain, "label")
stopifnot(all.equal(labels2, 1 - labels))
```

---

set\_field

*Set one attribute of a lgb.Dataset object*

---

### Description

Set one attribute of a lgb.Dataset

### Usage

```
set_field(dataset, field_name, data)

## S3 method for class 'lgb.Dataset'
set_field(dataset, field_name, data)
```

### Arguments

<code>dataset</code>	Object of class lgb.Dataset
<code>field_name</code>	String with the name of the attribute to set. One of the following. <ul style="list-style-type: none"> <li>• <code>label</code>: label lightgbm learns from ;</li> <li>• <code>weight</code>: to do a weight rescale ;</li> </ul>

- `group`: used for learning-to-rank tasks. An integer vector describing how to group rows together as ordered results from the same set of candidate results to be ranked. For example, if you have a 100-document dataset with `group = c(10, 20, 40, 10, 10, 10)`, that means that you have 6 groups, where the first 10 records are in the first group, records 11-30 are in the second group, etc.
- `init_score`: initial score is the base prediction lightgbm will boost from.

`data` The data for the field. See examples.

### Value

The `lgb.Dataset` you passed in.

### Examples

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)
lgb.Dataset.construct(dtrain)

labels <- lightgbm::get_field(dtrain, "label")
lightgbm::set_field(dtrain, "label", 1 - labels)

labels2 <- lightgbm::get_field(dtrain, "label")
stopifnot(all.equal(labels2, 1 - labels))
```

---

slice

*Slice a dataset*

---

### Description

Get a new `lgb.Dataset` containing the specified rows of original `lgb.Dataset` object

### Usage

```
slice(dataset, ...)

## S3 method for class 'lgb.Dataset'
slice(dataset, idxset, ...)
```

### Arguments

<code>dataset</code>	Object of class <code>lgb.Dataset</code>
<code>...</code>	other parameters (currently not used)
<code>idxset</code>	an integer vector of indices of rows needed

**Value**

constructed sub dataset

**Examples**

```
data(agaricus.train, package = "lightgbm")
train <- agaricus.train
dtrain <- lgb.Dataset(train$data, label = train$label)

dsub <- lightgbm::slice(dtrain, seq_len(42L))
lgb.Dataset.construct(dsub)
labels <- lightgbm::get_field(dsub, "label")
```

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